Quantitation Report (Qedit)

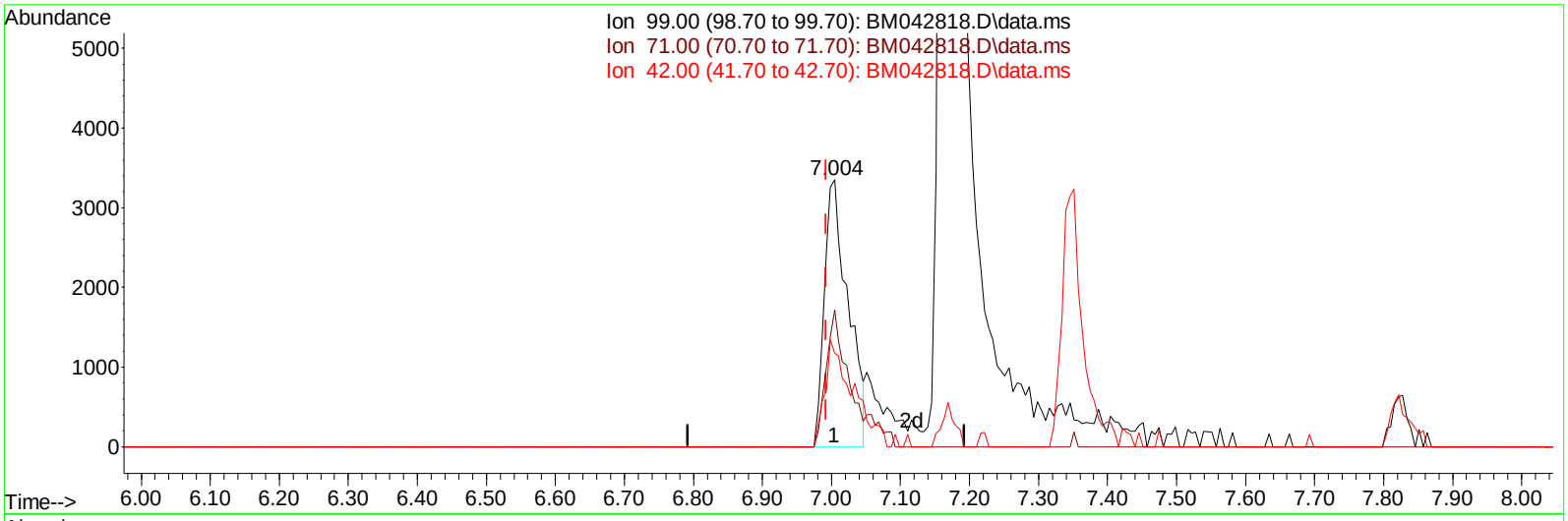
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111323\
Data File : BM042818.D
Acq On : 16 Nov 2023 22:45
Operator : MA/JU
Sample : 05268-09
Misc :
ALS Vial : 60 Sample Multiplier: 1

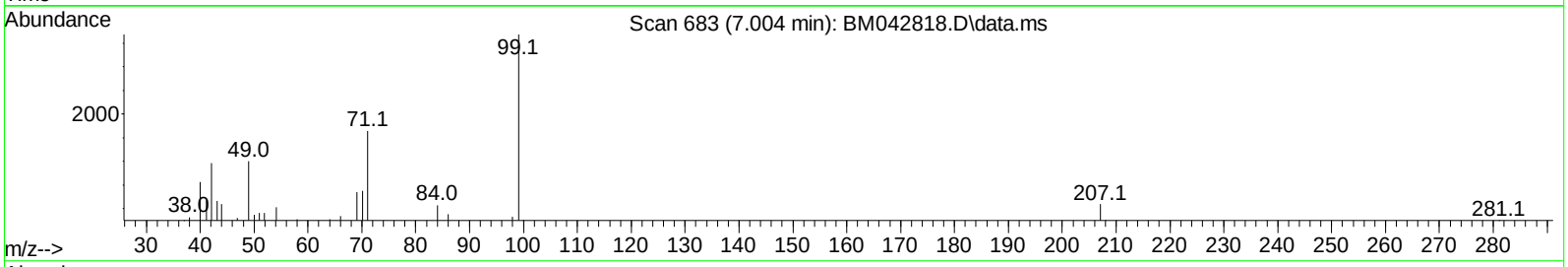
Quant Time: Nov 16 23:56:39 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 15 22:28:07 2023
Response via : Initial Calibration

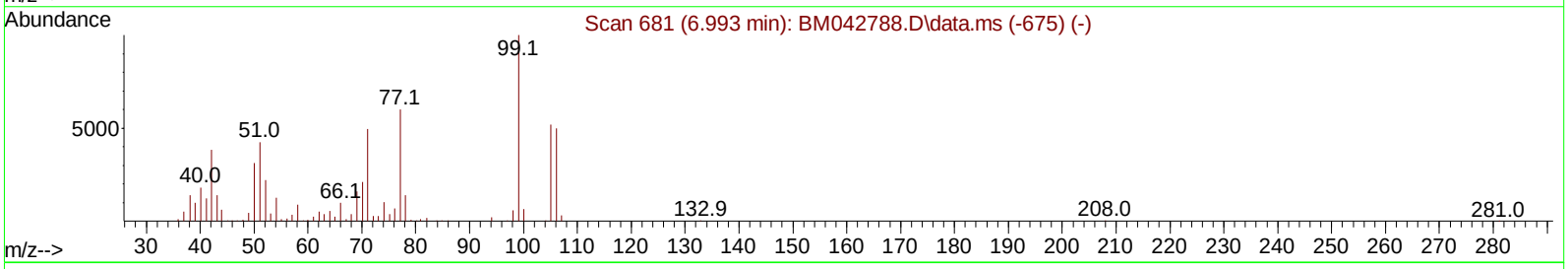
Instrument :
BNA_M
ClientSampleId :
DCKD1

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/17/2023
Supervised By :mohammad ahmed 11/19/2023







TIC: BM042818.D\data.ms

(7) Phenol-d5 (S)

7.004min (+ 0.012) 3.70 ng/ul

response 7984

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	43.50	51.18
42.00	32.20	35.00
0.00	0.00	0.00

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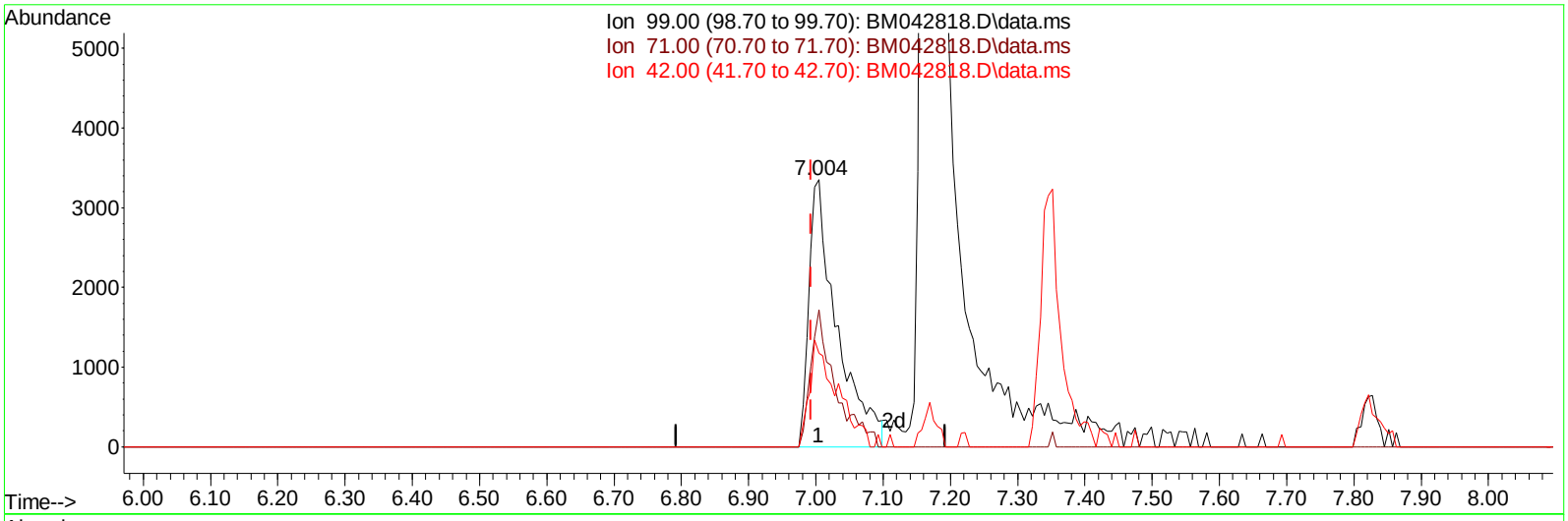

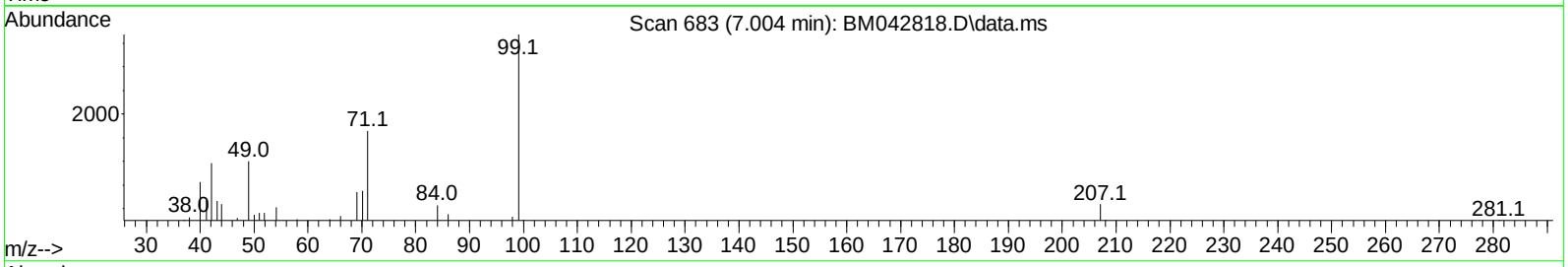

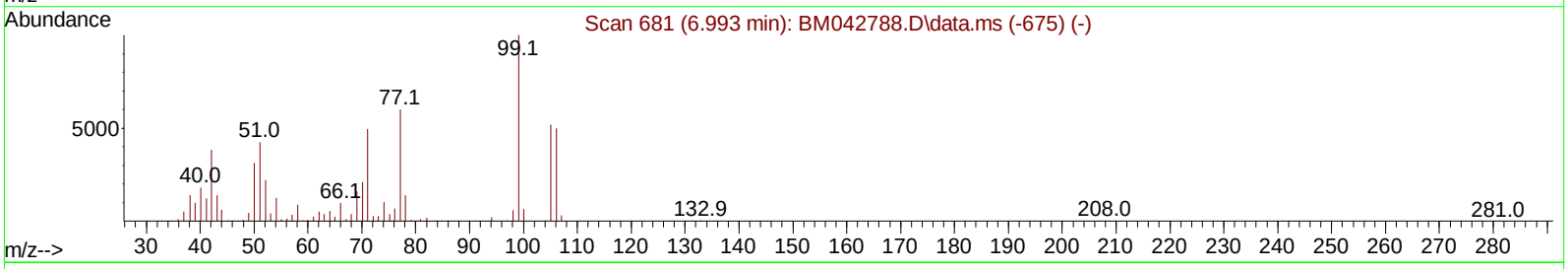

TIC: BM042818.D\data.ms

(7) Phenol-d5 (S)

7.004min (+ 0.012) 4.49 ng/ul m

response 9700

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	43.50	51.18
42.00	32.20	35.00
0.00	0.00	0.00

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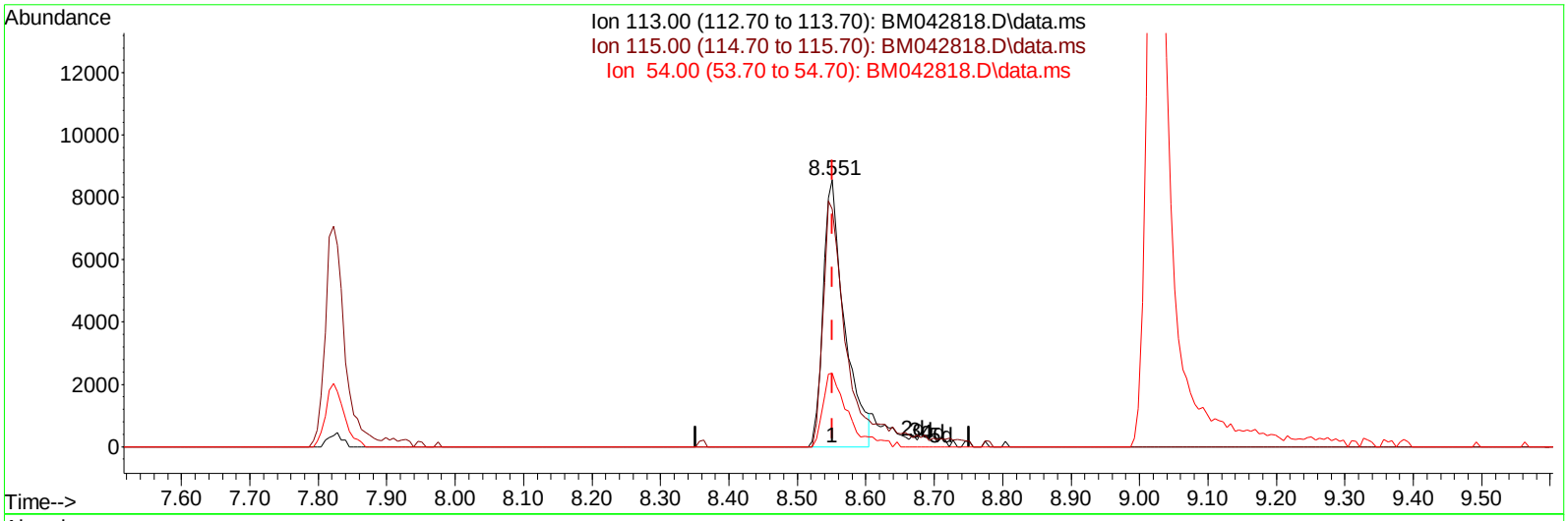

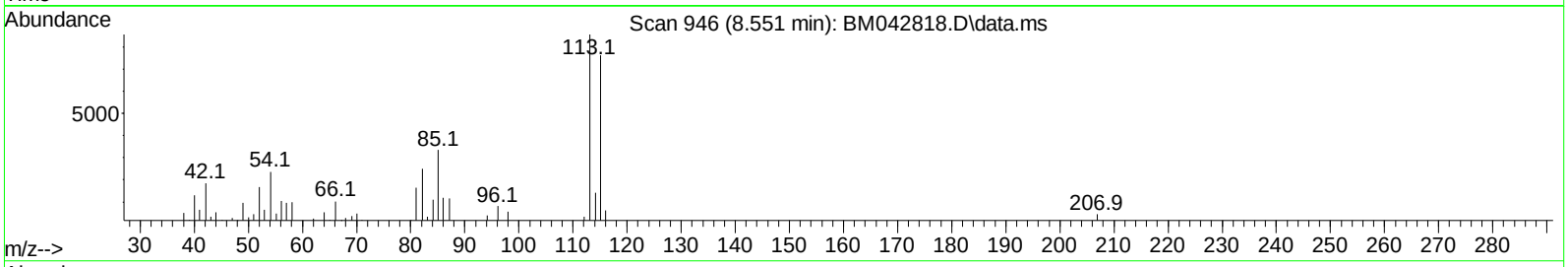

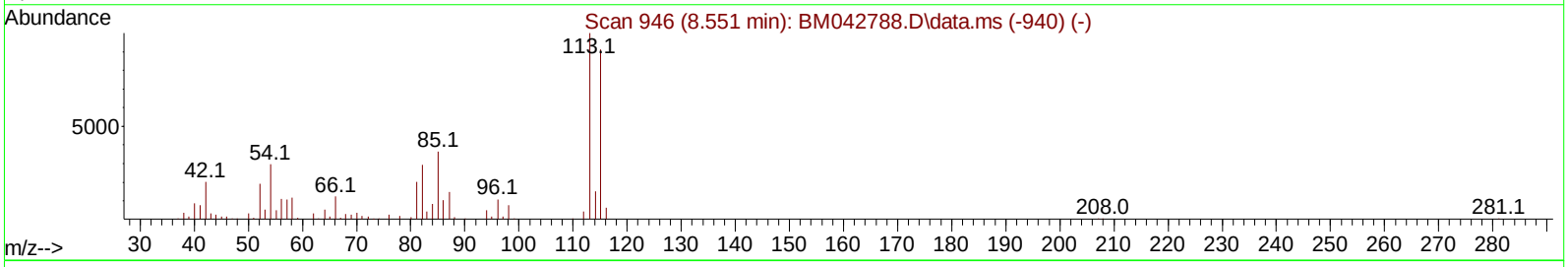

TIC: BM042818.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.551min (-0.000) 10.89 ng/ul

response 18517

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	91.70	88.91
54.00	24.00	27.59
0.00	0.00	0.00

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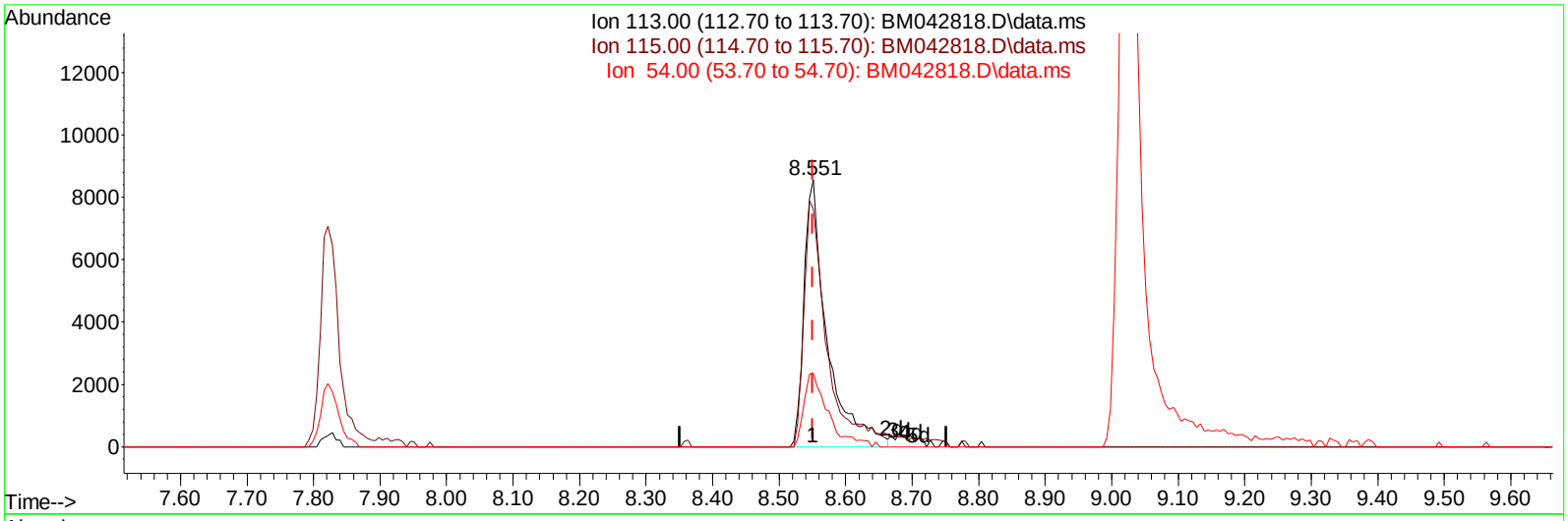

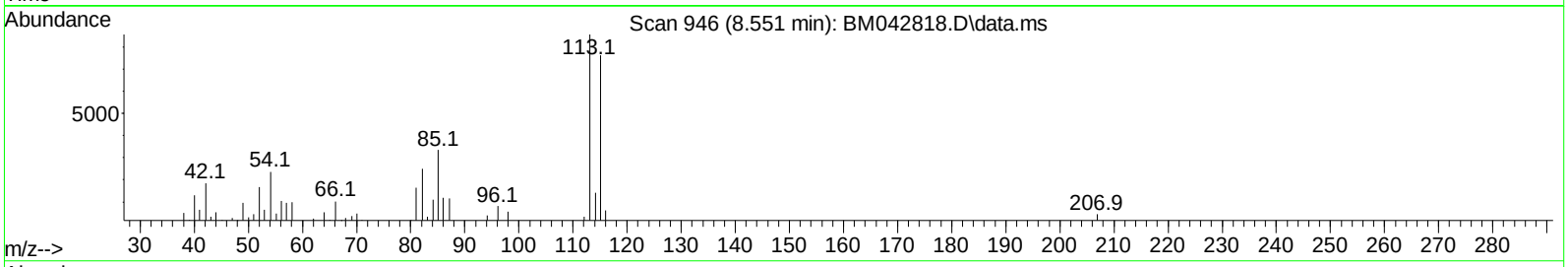

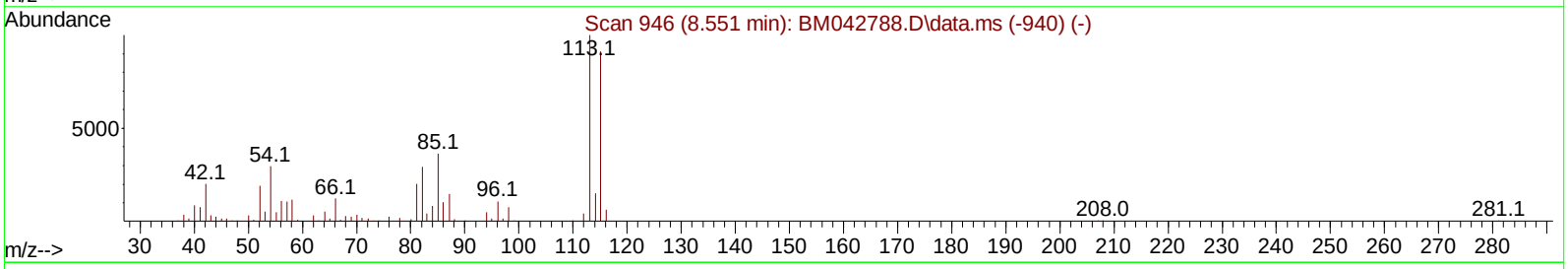

TIC: BM042818.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.551min (-0.000) 12.07 ng/ul m

response 20522

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	91.70	88.91
54.00	24.00	27.59
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	22196	20.000	ng/ul	0.00
20) Naphthalene-d8	10.634	136	89604	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.474	164	57886	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.233	188	124136	20.000	ng/ul	0.00
79) Chrysene-d12	21.415	240	127422	20.000	ng/ul	# 0.00
88) Perylene-d12	23.774	264	162233	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	1911	2.701	ng/uL	0.00
4) Pyridine-d5	3.628	84	10289	5.847	ng/ul	0.01
7) Phenol-d5	7.004	99	9700m	4.490	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.169	67	55028	36.252	ng/ul	0.00
11) 2-Chlorophenol-d4	7.345	132	37226	22.333	ng/ul	0.00
15) 4-Methylphenol-d8	8.551	113	20522m	12.065	ng/ul	0.00
21) Nitrobenzene-d5	9.028	128	23267	29.617	ng/ul	0.00
24) 2-Nitrophenol-d4	9.734	143	24105	26.229	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.263	165	38812	22.713	ng/ul	0.00
31) 4-Chloroaniline-d4	10.816	131	35649	16.431	ng/ul	0.00
46) Dimethylphthalate-d6	13.904	166	198017	36.144	ng/ul	0.00
49) Acenaphthylene-d8	14.180	160	204742	34.928	ng/ul	0.00
54) 4-Nitrophenol-d4	14.786	143	161	0.260	ng/ul	0.05
60) Fluorene-d10	15.474	176	172872	38.106	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.621	200	23024	24.486	ng/ul	0.00
73) Anthracene-d10	17.333	188	286495	41.181	ng/ul	0.00
81) Pyrene-d10	19.627	212	395268	47.371	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.621	264	403955	41.627	ng/ul	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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